

```
Data Path : Z:\HPCHEM1\BNA M\Data\BM031317\  
Data File : BM009223.D  
Acq On : 13 Mar 2017 18:22  
Operator : SJ/MA  
Sample : SSTDICV020  
Misc :  
ALS Vial : 8 Sample Multiplier: 1
```

Quant Time: Mar 13 18:56:31 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-2016\SOM-EPA-BM031317MA M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Mar 13 18:31:21 2017
Response via : Initial Calibration

Instrument :

BNA_M

ClientSampleId :

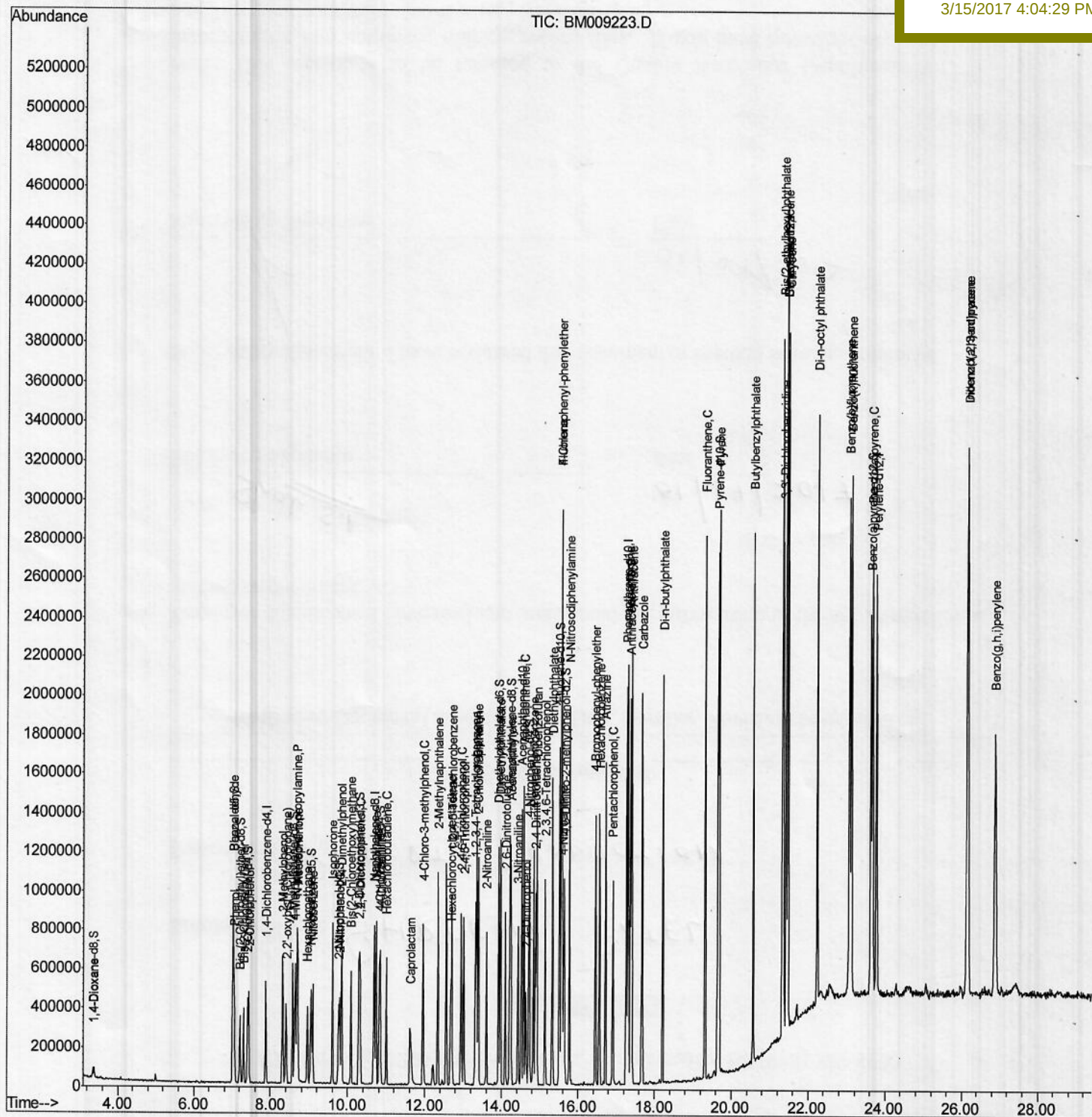
SICV66

Manual Integrations

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Quantitation Report (Qedit)

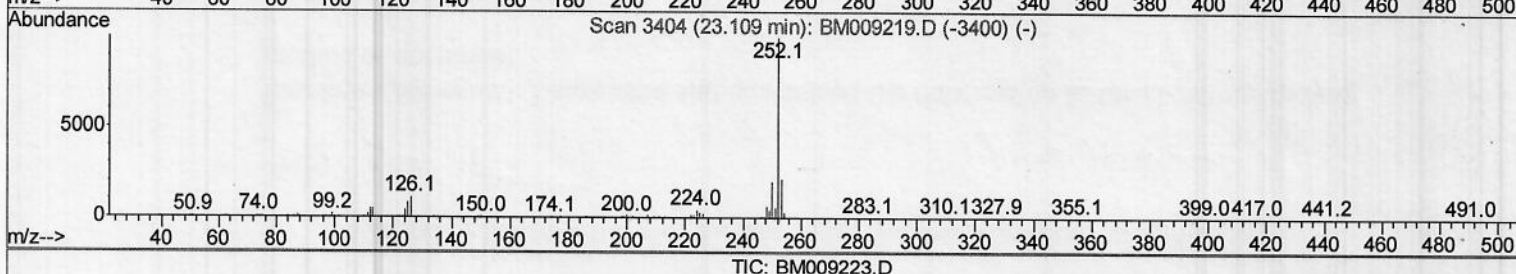
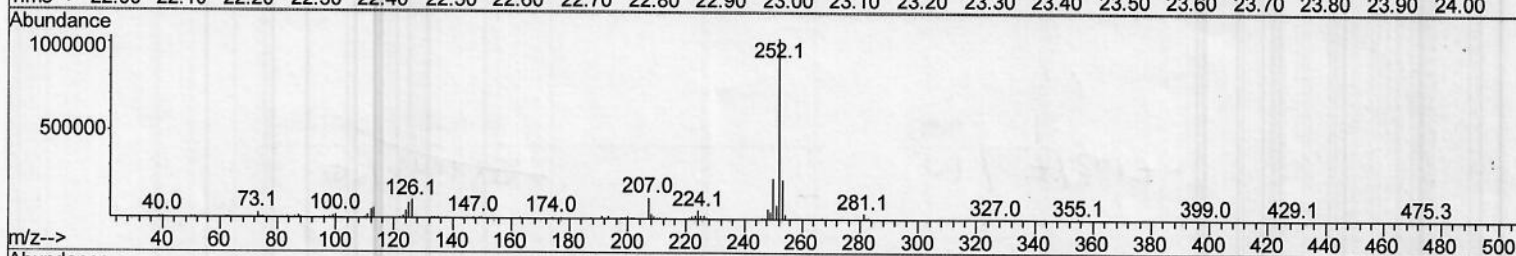
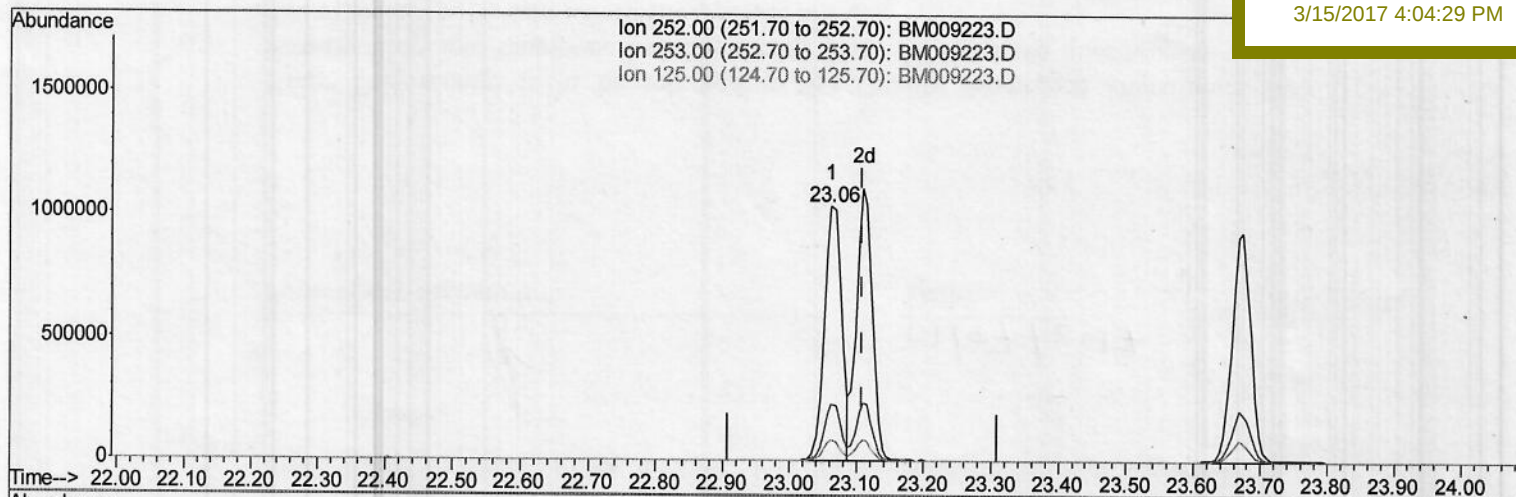
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TIC: BM009223.D

(88) Benzo(k)fluoranthene

23.062min (-0.047) 20.61ng/ul

response 1864178

Ion	Exp%	Act%
252.00	100	100
253.00	21.60	22.16
125.00	10.20	8.21
0.00	0.00	0.00

Quantitation Report (Qedit)

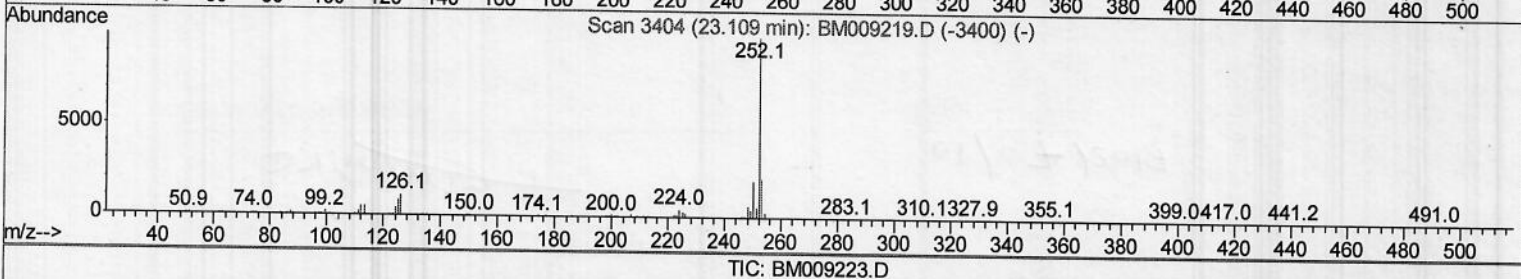
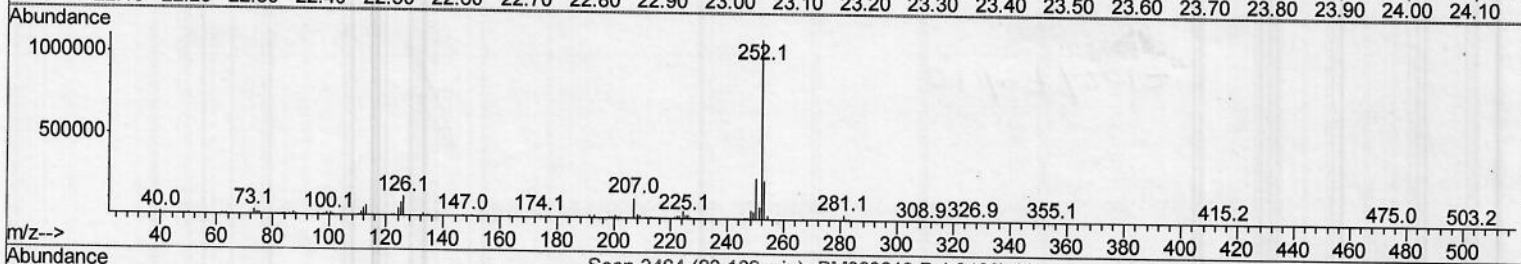
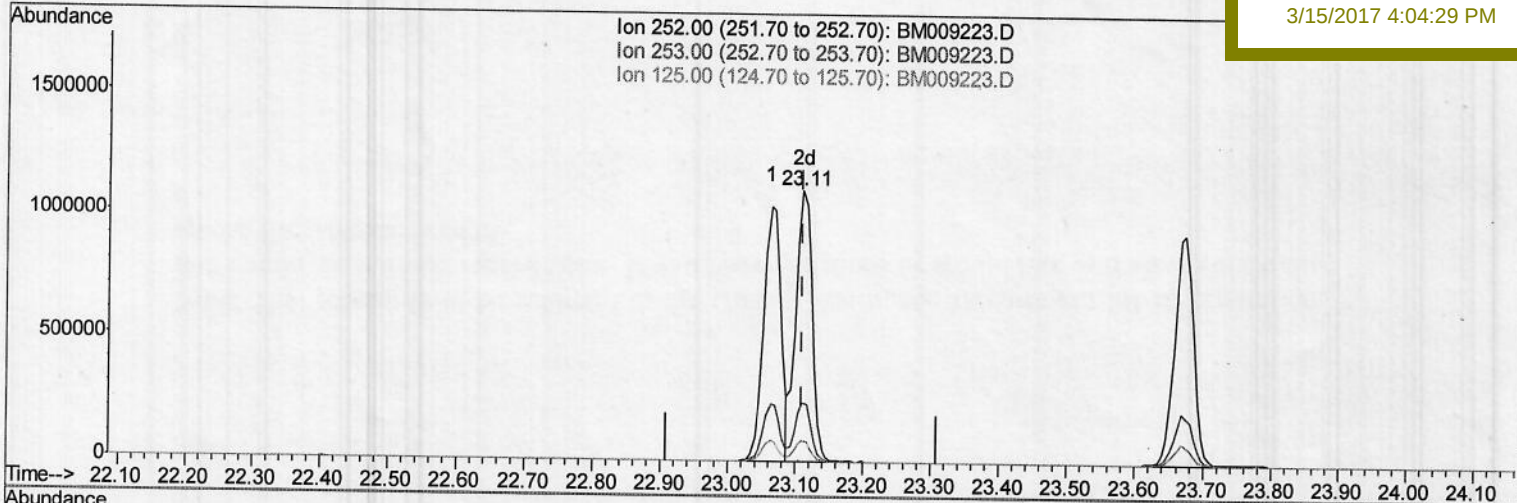
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 Client Sampled :
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Manual Integrations
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(88) Benzo(k)fluoranthene

23.109min (+0.000) 20.30ng/ul m

response 1836118

Ion	Exp%	Act%
252.00	100	100
253.00	21.60	21.15
125.00	10.20	7.77#
0.00	0.00	0.00

50
 03/18/2017

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	119124	20.00	nq/ul	0.00
18) Naphthalene-d8	10.68	136	572226	20.00	nq/ul	0.00
35) Acenaphthene-d10	14.52	164	423943	20.00	nq/ul	0.00
61) Phenanthrene-d10	17.27	188	1128562	20.00	nq/ul	0.00
77) Chrysene-d12	21.44	240	1600684	20.00	nq/ul	0.00
85) Perylene-d12	23.77	264	1600989	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.35	96	22024	8.14	nq/uL	0.00
5) Phenol-d5	7.03	99	234359	18.19	nq/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.21	67	162229	19.33	nq/ul	0.00
9) 2-Chlorophenol-d4	7.41	132	149990	19.05	nq/ul	0.00
13) 4-Methylphenol-d8	8.58	113	192126	18.69	nq/ul	0.00
19) Nitrobenzene-d5	9.05	128	84464	20.67	nq/ul	0.00
22) 2-Nitrophenol-d4	9.76	143	91326	20.13	nq/ul	0.00
26) 2,4-Dichlorophenol-d3	10.29	165	195611	20.19	nq/ul	0.00
29) 4-Chloroaniline-d4	10.82	131	237362	20.54	nq/ul	0.00
43) Dimethylphthalate-d6	13.92	166	720637	22.38	nq/ul	0.00
46) Acenaphthylene-d8	14.22	160	819345	22.13	nq/ul	0.00
51) 4-Nitrophenol-d4	14.71	143	135444	21.19	nq/ul	0.00
57) Fluorene-d10	15.51	176	628821	21.89	nq/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.63	200	121908	18.06	nq/ul	0.00
70) Anthracene-d10	17.36	188	1053737	19.73	nq/ul	0.00
78) Pyrene-d10	19.66	212	1307307	20.23	nq/ul	0.00
89) Benzo(a)pyrene-d12	23.62	264	1454302	19.86	ng/ul	0.00

Target Compounds

					Ovalue
2) 1,4-Dioxane	3.38	88	24952	8.70	nq/uL# 75
4) Benzaldehyd	7.03	77	192892	22.33	nq/ul# 76
6) Phenol	7.06	94	248493	18.06	nq/ul# 79
8) Bis(2-Chloroethyl) ether	7.30	93	185173	18.79	nq/ul# 74
10) 2-Chlorophenol	7.44	128	155820	19.19	nq/ul# 75
11) 2-Methylphenol	8.31	108	178170	17.87	nq/ul 95
12) 2,2'-oxybis(1-Chloropropan	8.41	45	311534	18.97	nq/ul# 86
14) Acetophenone	8.70	105	322103	18.65	nq/ul# 77
15) N-Nitroso-di-n-propylamine	8.68	70	202029	19.08	nq/ul# 85
16) 4-Methylphenol	8.64	108	206681	18.29	nq/ul 94
17) Hexachloroethane	8.96	117	67930	19.10	nq/ul 98
20) Nitrobenzene	9.09	77	272138	19.97	nq/ul# 86
21) Isophorone	9.60	82	532430	19.75	nq/ul# 93
23) 2-Nitrophenol	9.79	139	100834	20.42	nq/ul# 72
24) 2,4-Dimethylphenol	9.85	107	254301	20.07	nq/ul 89
25) Bis(2-Chloroethoxy)methane	10.09	93	293915	19.89	nq/ul 98
27) 2,4-Dichlorophenol	10.32	162	186674	19.46	nq/ul 95
28) Naphthalene	10.73	128	573085	19.53	nq/ul 98
30) 4-Chloroaniline	10.85	127	243403	19.88	nq/ul 99
31) Hexachlorobutadiene	11.00	225	131556	19.90	nq/ul 98
32) Caprolactam	11.61	113	76916	19.20	nq/ul# 58
33) 4-Chloro-3-methylphenol	11.96	107	254143	20.19	nq/ul 93
34) 2-Methylnaphthalene	12.34	142	455592	19.50	ng/ul 96

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36) 1,2,4,5-Tetrachlorobenzene	12.70	216	279043	21.88	nq/ul	97
37) Hexachlorocyclopentadiene	12.67	237	145214	20.14	nq/ul	96
38) 2,4,6-Trichlorophenol	12.95	196	176628	21.85	nq/ul	95
39) 2,4,5-Trichlorophenol	13.02	196	203420	21.96	nq/ul	99
40) 1,1'-Biphenyl	13.35	154	649243	21.86	nq/ul	94
41) 2-Chloronaphthalene	13.39	162	496783	21.61	nq/ul	96
42) 2-Nitroaniline	13.60	65	213631	22.63	nq/ul#	71
44) Dimethylphthalate	13.97	163	722390	22.37	nq/ul#	99
45) 2,6-Dinitrotoluene	14.10	165	135671	21.84	nq/ul#	69
47) Acenaphthylene	14.24	152	818416	22.08	nq/ul	98
48) 3-Nitroaniline	14.43	138	148057	22.38	nq/ul	82
49) Acenaphthene	14.59	153	568798	21.81	nq/ul	96
50) 2,4-Dinitrophenol	14.63	184	79486	19.52	nq/ul#	76
52) 4-Nitrophenol	14.72	109	154260	20.66	nq/ul#	76
53) Dibenzofuran	14.92	168	841196	21.89	nq/ul	98
54) 2,4-Dinitrotoluene	14.89	165	216344	22.60	nq/ul#	76
55) 2,3,4,6-Tetrachlorophenol	15.14	232	195560	22.00	nq/ul	93
56) Diethylphthalate	15.33	149	745703	22.05	nq/ul	99
58) Fluorene	15.57	166	716649	21.61	nq/ul	98
59) 4-Chlorophenyl-phenylether	15.56	204	373992	21.94	nq/ul	88
60) 4-Nitroaniline	15.59	138	168733	22.46	nq/ul#	66
63) 4,6-Dinitro-2-methylphenol	15.65	198	137302	19.36	nq/ul	93
64) N-Nitrosodiphenylamine	15.77	169	625552	19.72	nq/ul	98
65) 4-Bromophenyl-phenylether	16.46	248	237903	19.51	nq/ul	89
66) Hexachlorobenzene	16.57	284	265923	19.61	nq/ul#	86
67) Atrazine	16.72	200	256229	19.32	nq/ul	94
68) Pentachlorophenol	16.91	266	156427	17.56	nq/ul	97
69) Phenanthrene	17.31	178	1233231	19.64	nq/ul	100
71) Anthracene	17.40	178	1266007	19.77	nq/ul	98
72) 1,2,3,4-Tetrachlorobenzene	13.32	216	260179	19.62	nq/uL	99
73) Pentachlorobenzene	14.83	250	284764	20.00	nq/uL	97
74) Carbazole	17.67	167	1160434	19.41	nq/ul	98
75) Di-n-butylphthalate	18.22	149	1327392	19.26	nq/ul#	98
76) Fluoranthene	19.32	202	1578822	19.94	nq/ul	97
79) Pvrene	19.68	202	1668600	20.06	nq/ul	98
80) Butylbenzylphthalate	20.57	149	654713	19.59	nq/ul#	80
81) 3,3'-Dichlorobenzidine	21.36	252	650109	21.44	nq/ul#	98
82) Benzo(a)anthracene	21.42	228	1825108	19.95	nq/ul	98
83) Bis(2-ethylhexyl)phthalate	21.33	149	987012	18.97	nq/ul#	99
84) Chrysene	21.47	228	1693073	19.62	nq/ul	99
86) Di-n-octyl phthalate	22.24	149	1719688	19.97	nq/ul#	89
87) Benzo(b)fluoranthene	23.06	252	1864178	19.70	nq/ul#	98
88) Benzo(k)fluoranthene	23.11	252	1836118m	20.30	nq/ul	98
90) Benzo(a)pyrene	23.67	252	1807425	19.65	nq/ul#	96
91) Indeno(1,2,3-cd)pyrene	26.16	276	2117855	19.19	nq/ul#	90
92) Dibenzo(a,h)anthracene	26.16	278	1783054	19.11	nq/ul#	94
93) Benzo(a,h,i)perylene	26.89	276	1783226	19.53	nq/ul#	90

(#) = qualifier out of range (m) = manual integration (+) = signals summed